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If more than one search is su ************************ Please provide a detailed statement of Include the elected species or structure utility of the invention. Define any ter known. Please attach a copy of the cov	the search topic, and descries, keywords, synonyms, ac	ibe as specifically as possible the sucronyms, and registry numbers, and	**************************************
Title of Invention: Ogan	ic Electrol	umin'eseent	Danie
Inventors (please provide full names	William R.	egley, Tukaran	Habus
Manju Rajeswa	ran, David (riesen Natasha	Andrieus le
Earliest Priority Filing Date:	11/4/2003		- Jimires ay
For Sequence Searches Only Please inc appropriate serial number.	/ clude all pertinent informatio	n (parent, child, divisional, or issued	patent numbers) along with the
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Date Completed: 10 - 20 - 1	Bibliographic	Dr.Link	
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PTO-1590 (8-01)		(

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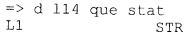
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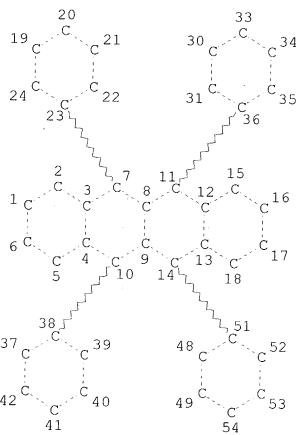
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L7 L8 L9 L10	FILE	'REGISTRY' ENTERED AT 15:50:58 ON 20 OCT 2004 8 S L1 185 S L1 FUL SAV L8 GAR916/A 0 S L5 SSS SAM SUB=L8 1 S L5 SSS FUL SUB=L8 SAV L10 GAR916A/A
L11	FILE	'CAOLD' ENTERED AT 15:54:41 ON 20 OCT 2004 0 S L10
L12	FILE	'ZCAPLUS' ENTERED AT 15:54:49 ON 20 OCT 2004 1 S L10
L13 L14	FILE	'REGISTRY' ENTERED AT 15:55:12 ON 20 OCT 2004 1 S L3 SSS SAM SUB=L8 2 S L3 SSS FUL SUB=L8 SAV L14 GAR241/A
L15	FILE	'CAOLD' ENTERED AT 15:56:19 ON 20 OCT 2004 1 S L14
L16	FILE	'ZCAPLUS' ENTERED AT 15:58:50 ON 20 OCT 2004 2 S L14
L17 L18	FILE	'REGISTRY' ENTERED AT 15:59:03 ON 20 OCT 2004 0 S L2 SSS SAM SUB=L8 1 S L2 SSS FUL SUB=L8 SAV L18 GAR040/A

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FILE 'ZCAPLUS' ENTERED AT 16:00:50 ON 20 OCT 2004 L20 1 S L18

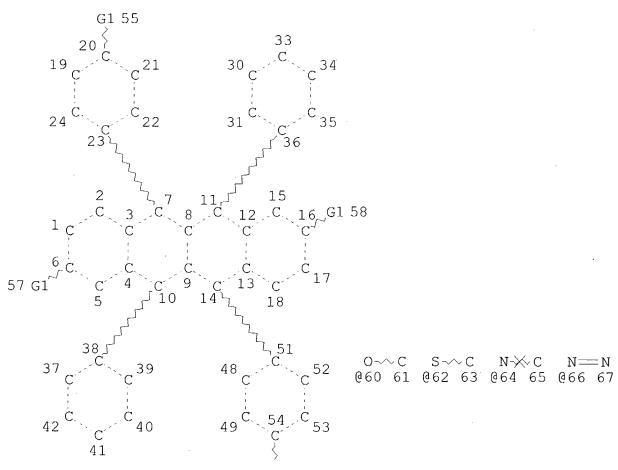
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NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RSPEC I NUMBER OF NODES IS 42 STEREO ATTRIBUTES: NONE L3 STR



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Page 2-A

VAR G1=60/62/64/66

NODE ATTRIBUTES:

NSPEC IS RC ΑT 61 NSPEC IS RC ΑT 63 NSPEC IS RC AT64 NSPEC IS RC AT65 CONNECT IS E2 RC AT DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 54

STEREO ATTRIBUTES: NONE

L8 185 SEA FILE=REGISTRY SSS FUL L1

L14 2 SEA FILE=REGISTRY SUB=L8 SSS FUL L3

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SEARCH TIME: 00.00.01

2 ANSWERS

=> file caold

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FILE LAST UPDATED: 01 May 1997 (19970501/UP)

=> d 115 1 all hitstr

L15 ANSWER 1 OF 1 CAOLD COPYRIGHT 2004 ACS on STN

AN CA51:12055c CAOLD

TI diarylphenylenenaphthacene derived from a tetramethoxyrubrene

AU Perronnet, Jacques

IT 118769-17-8 119504-35-7 121544-89-6

IT 118769-17-8

RN 118769-17-8 CAOLD

CN Naphthacene, 2,8-dimethoxy-5,11-bis(p-methoxyphenyl)-6,12-diphenyl-(6CI) (CA INDEX NAME)

=> file zcaplus FILE 'ZCAPLUS' ENTERED AT 16:05:25 ON 20 OCT 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

=> d l16 1-2 ibib abs hitstr hitrn

L16 ANSWER 1 OF 2 ZCAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 1956:16251 ZCAPLUS

DOCUMENT NUMBER: 50:16251

ORIGINAL REFERENCE NO.: 50:3369c-e

TITLE: Determination of the structures of eight methoxy tetraphenylnaphthacenes

AUTHOR(S): Dufraisse, Charles; Etienne, Andre; Valls, Jaime

SOURCE: Compt. rend. (1955), 240, 2097-2100

DOCUMENT TYPE: Journal LANGUAGE: Unavailable

AB cf. C.A. 49, 6896h. The structures of the following substituted naphthacenes were assigned: 2,8-(MeO)2, 5,6,11,12-Ph4 (I), m. 259.degree.; 2-MeO, 11-(p-MeOC6H4), 5,6,12-Ph3 (II), m. 227.degree.; 5,11-(p-MeOC6H4)2, 6,12-Ph2 (III), m. 252.degree.; 2,8-(MeO)2, 5,11-(p-MeOC6H4)2, 6,12-Ph2 (IV), m. 257.degree.; 2,8-(MeO)2, 6,12-(p-MeOC6H4)2, 5,11-Ph2 (V), m. 250.degree.; 8-MeO,

5,6,12-(p-MeOC6H4)3, 11-Ph (VI), m. 226.degree.; 5,6,11,12-(p-MeOC6H4)4 (VII), m. 282.degree.; and 2,8-(MeO)2, 5,6,11,12-(p-MeOC6H4)4 (VIII), m. 255.degree. The structures of III, IV, and VIII are known; those of the other compds. are assigned on the basis of m.ps., relative adsorption on Al2O3 on chromatog. (in order of decreasing adsorption, III, II, I; and VII, VI, V), and UV absorption spectra (the appearance of new absorption peaks in the spectra of I, II, IV, V, VI, and VIII is attributed to direct attachment of MeO to the naphthacene ring).

118769-17-8, Naphthacene, 2,8-dimethoxy-5,11-bis(p-methoxyphenyl)-6,12-diphenyl-694489-88-8, Naphthacene, 2,8-dimethoxy-5,6,11,12-tetrakis-(p-methoxyphenyl)-(prepn. of)

RN 118769-17-8 ZCAPLUS

CN Naphthacene, 2,8-dimethoxy-5,11-bis(p-methoxyphenyl)-6,12-diphenyl-(6CI) (CA INDEX NAME)

RN 694489-88-8 ZCAPLUS

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118769-17-8, Naphthacene, 2,8-dimethoxy-5,11-bis(p-ΙT methoxyphenyl)-6,12-diphenyl- 694489-88-8, Naphthacene, 2,8-dimethoxy-5,6,11,12-tetrakis-(p-methoxyphenyl)-(prepn. of)

ANSWER 2 OF 2 ZCAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER:

1955:35916 ZCAPLUS

DOCUMENT NUMBER:

49:35916

ORIGINAL REFERENCE NO.:

49:6896h-i,6897a-b

TITLE:

Systematic enumeration of the methoxylated

AUTHOR(S):

rubrenes obtainable by rubrenic reactions

SOURCE:

Dufraisse, Charles; Etienne, Andre; Valls, Jaime Compt. rend. (1954), 239, 1101-4

DOCUMENT TYPE:

Journal

LANGUAGE:

Unavailable

Eight methoxylated rubrenes were obtained from the 5 p-methoxy triphenylpropargyl alcs. by rubrenic condensation of the hydrochlorides of the corresponding propargyl alcs. or of their chloride esters. This is exactly the no. of methoxylated rubrenes predicted on the basis of a centrosym. reaction scheme. rubrenes obtained were: 9,11-di(p-methoxyphenyl)-10,12diphenylnaphthacene [from Ph2C(OH)C .tplbond. CC6H4OMe-p or p-MeOC6H4PhC(OH)C.tplbond.CPh]; 9-p-methoxyphenyl-10,11,12-triphenyl-

2-methoxynaphthacene, m. 227.degree., and 9,10,11,12-tetraphenyl-2,6dimethoxynaphthacene, m. 259.degree. [from p-MeOC6H4PhC(OH)C.tplbond.CPh]; 9,11-di(p-methoxyphenyl)-10,12diphenyl-2,6-dimethoxynaphthacene, m. 257.degree. [from (p-MeOC6H4) 2C (OH) C.tplbond.CPh]; 9,11-di(p-methoxyphenyl)-10,12diphenyl-3,7-dimethoxynaphthacene, m. 250.degree.; 9,10,11-tri(p-methoxyphenyl)-12-phenyl-7-methoxynaphthacene, m. 226.degree. (C6H6 solvate, m. 195.degree.); and 9,10,11,12-tetra(pmethoxyphenyl) naphthacene, m. 282.degree. (C6H6 solvate, m. 245.degree.) [from p-MeOC6H4PhC(OH)C.tplbond.CC6H4OMe-p]; and 9,10,11,12-tetra(p-methoxyphenyl)-2,6-dimethoxynaphthacene, m. 255.degree. [from (p-MeOC6H4)2C(OH)C.tplbond.CC6H4OMe-p]. Addnl. products, as yet unidentified, were also obtained; these were, however, shown not to have the rubrenic structure. 118769-17-8, Naphthacene, 2,8-dimethoxy-5,11-bis(pmethoxyphenyl)-6,12-diphenyl- 694489-88-8, Naphthacene, 2,8-dimethoxy-5,6,11,12-tetrakis-(p-methoxyphenyl)-(prepn. of)

RN 118769-17-8 ZCAPLUS

ΙT

CN Naphthacene, 2,8-dimethoxy-5,11-bis(p-methoxyphenyl)-6,12-diphenyl-(6CI) (CA INDEX NAME)

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